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AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions of the claims and listing of the claims in the application:

1. (Currently Amended) A substituted tetracycline compound, wherein said compound is of the formula:

(I)

wherein:

X is CHC(R¹³Y'Y), CR⁶'R⁶, S, NR⁶, or O:

R² is hydrogen, alkyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic or a prodrug moiety;

R⁴ and R⁴ are each hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic or a prodrug moiety;

R², R³, R¹⁰, R¹¹ and R¹² are each hydrogen or a pro-drug moiety;

R⁵ is hydrogen, hydroxyl, or a prodrug moiety:

R⁶, R⁶, and R⁸ are each independently hydrogen, alkyl, alkenyl, alkynyl, aryl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, or halogen;

R⁷ is hydrogen, dialkylamino, or NR^{7c}C(=W')WR^{7a};

R⁸ is hydrogen:

R¹³ is hydrogen, hydroxy, alkyl; alkenyl; alkynyl; alkoxy; alkylthio; alkylsulfinyl; alkylsulfonyl; alkylamino; or an arylalkyl;

Y' and Y are each independently hydrogen; halogen; hydroxyl; cyano, sulfhydryl; amino; alkyl; alkenyl; alkynyl; alkoxy; alkylthio; alkylsulfinyl; alkylsulfonyl; alkylamino; or an arylalkyl;

R⁹ is hydrogen, or NR^{9c}C(=Z')ZR^{9a};

Z is O;

Z' is O or S;

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 R^{9a} is unsubstituted C_3 - C_{10} alkyl, substituted alkyl, substituted or unsubstituted alkenyl, substituted or unsubstituted alkynyl, substituted or unsubstituted alkoxy, substituted or unsubstituted alkylthio, substituted or unsubstituted alkylsulfinyl, substituted or unsubstituted alkylsulfonyl, substituted or unsubstituted arylsulfonyl, substituted or unsubstituted alkoxycarbonyl, substituted or unsubstituted arylcarbonyl, substituted or unsubstituted alkylamino, substituted or unsubstituted arylalkyl, substituted or unsubstituted aryl, substituted or unsubstituted heterocyclic, substituted or unsubstituted heteroaromatic, a steroid, absent, or a prodrug moiety, wherein said substituted alkyl is substituted with halogen, amino, hydroxyl, alkoxy, alkylcarbonyloxy, alkyloxycarbonyl, arylcarbonyloxy, alkoxycarbonylamino, alkoxycarbonyloxy, aryloxycarbonyloxy, carboxylate, alkylcarbonyl, alkylaminoacarbonyl, arylalkyl aminocarbonyl, alkenylaminocarbonyl, alkylcarbonyl, arylcarbonyl, aminoalkyl, arylalkylcarbonyl, alkenylcarbonyl, alkoxycarbonyl, silyl, aminocarbonyl, alkylthiocarbonyl, phosphate, aralkyl, phosphonato, phosphinato, cyano, acylamino, amido, imino, sulfhydryl, alkylthio, sulfate, arylthio, thiocarboxylate, alkylsulfinyl, sulfonato, sulfamoyl, sulfonamido, nitro, cyano, azido, heterocyclyl, alkylaryl, aryl or heteroaryl; further wherein said substituted alkenyl, substituted alkynyl, substituted alkoxy, substituted alkylthio, substituted alkylsulfinyl, substituted alkylsulfonyl, substituted arylsulfonyl, substituted alkoxycarbonyl, substituted arylcarbonyl, substituted alkylamino, substituted arylalkyl, substituted aryl, substituted heterocyclic, or substituted heteroaromatic is substituted with halogen, amino, alkyl, alkenyl, alkynyl, hydroxyl, alkoxy, alkylcarbonyloxy, alkyloxycarbonyl, arylcarbonyloxy, alkoxycarbonylamino, alkoxycarbonyloxy, aryloxycarbonyloxy, carboxylate, alkylcarbonyl, alkylaminoacarbonyl, arylalkyl aminocarbonyl, alkenylaminocarbonyl, alkylcarbonyl, arylcarbonyl, aminoalkyl, arylalkylcarbonyl, alkenylcarbonyl, alkoxycarbonyl, silyl, aminocarbonyl, alkylthiocarbonyl, phosphate, aralkyl, phosphonato, phosphinato, cyano, acylamino, amido, imino, sulfhydryl, alkylthio, sulfate, arylthio, thiocarboxylate, alkylsulfinyl, sulfonato, sulfamoyl, sulfonamido, nitro, cyano, azido, heterocyclyl, alkylaryl, aryl or heteroaryl;

R^{9c} is hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, arylsulfonyl, alkoxycarbonyl, arylcarbonyl, alkylamino, arylalkyl, aryl, heterocyclic or heteroaromatic;

W is CR^{7d}R^{7e}, NR^{7b} or O;

W' is O or S; and

R^{7a}, R^{7b}, R^{7c}, R^{7d}, and R^{7e} are each independently hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, arylsulfonyl, alkoxycarbonyl, arylcarbonyl, alkylamino, arylalkyl,

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aryl, heterocyclic, heteroaromatic, absent, or a prodrug moiety, and R^{7d} and R^{7e} may be linked to form a ring;

and pharmaceutically acceptable salts thereof, provided that at least one of R⁹ is not hydrogen when R⁷ is hydrogen or dialkylamino.

- The compound of claim 1, wherein R², R², R³, R⁸, R¹⁰, R¹¹, and R¹² are each 2. (Original) hydrogen.
- The compound of claim 2, wherein R⁴ and R⁴ are each alkyl. 3. (Original)
- The compound of claim 3, wherein R⁴ and R⁴ are each methyl 4. (Original)
- 5. (Cancelled)
- The compound of claim 4, wherein R⁵ is hydrogen. (Original) 6.
- The compound of claim 6, wherein X is CH₂, and R⁷ is hydrogen. 7. (Original)
- The compound of claim 1406, wherein X is CH₂, and R⁷ is 8. (Currently Amended) $N(Me)_2$.
- The compound of claim 4, wherein R⁵ is hydroxyl or a prodrug moiety, and X 9. (Original) is CHR⁶.
- The compound of claim 9, wherein R⁵ is hydroxyl and R⁶ is CH₃. 10. (Original)
- The compound of claim 1, wherein R⁹ is NR^{9c}C(=Z')ZR^{9a}. 11. (Original)
- The compound of claim 11, wherein R^{9c} is hydrogen. 12. (Original)
- 13. (Original) The compound of claim 11, wherein Z' is oxygen.
- 14. (Original) The compound of claim 11, wherein Z' is sulfur.

- 15. (Cancelled)
- 16. (Original) The compound of claim 13 or 14, wherein Z is oxygen.
- 17. (Cancelled)
- 18. (Cancelled)
- 19. **(Previously presented)** The compound of claim 11, wherein R^{9a} is selected from the group consisting of substituted alkyl, alkynyl, aryl, arylalkyl, or heteroaromatic.
- 20. (Cancelled)
- 21. (Original) The compound of claim 19, wherein said substituted alkyl is substituted with one or more substituents selected from the group consisting of alkoxycarbonyl, amino, arylcarbonyl, halogen, hydroxy, alkylamino, alkoxy, or aryl.
- 22. (Cancelled)
- 23. **(Previously presented)** The compound of claim 19, wherein said substituted alkyl is substituted with an aryl group.
- 24. (Original) The compound of claim 23, wherein said aryl group is phenyl.
- 25. **(Previously presented)** The compound of claim 19, wherein said said substituted alkyl is substituted with one or more halogens.
- 26. (Original) The compound of claim 24, wherein said halogen is bromine.
- 27.-29 (Cancelled)
- 30. (Original) The compound of claim 19, wherein R^{9a} is substituted or unsubstituted aryl.

31. (Original) The compound of claim 30, wherein said substituted or unsubstituted aryl is naphthyl.

32. (Original) The compound of claim 30, wherein said substituted or unsubstituted aryl is of the formula:

- 33. (Original) The compound of claim 30, wherein said substituted or unsubstituted aryl is phenyl.
- 34. (Original) The compound of claim 33, wherein said aryl is substituted with one or more substituents selected from the group consisting of alkyl, alkenyl, alkynyl, aryl, alkoxy, aryloxy, alkylcarbonyl, arylcarbonyl, alkoxycarbonyl, aryloxycarbonyl, amido, halogen, nitro, azo, alkyl sulfonyl, and arylsulfonyl.
- 35. (Original) The compound of claim 34, wherein said substituent is alkyl.
- 36. (Original) The compound of claim 35, wherein said alkyl is unsubstituted.
- 37. (Original) The compound of claim 35, wherein said alkyl is methyl.
- 38. (Original) The compound of claim 35, wherein said alkyl is substituted with one or more halogens.
- 39. (Original) The compound of claim 34, wherein said substituent is methoxy.

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40. **(Original)** The compound of claim 34, wherein said substituent is selected from the group consisting of alkylcarbonyl, arylcarbonyl, alkoxycarbonyl, aryloxycarbonyl, and amido.

41-55. (Cancelled)

- 56. (Original) The compound of claim 1, wherein R⁷ is NR^{7c}C(=W')WR^{7a}.
- 57. (Original) The compound of claim 56, wherein R⁹ is hydrogen.
- 58. (Original) The compound of claim 57, wherein R^{7c} is hydrogen.
- 59. (Original) The compound of claim 57, wherein W' is oxygen.
- 60. (Original) The compound of claim 57, wherein W' is sulfur
- 61. (Original) The compound of claims 59 or 60, wherein W is NR^{7b}.
- 62. (Original) The compound of claims 59 or 60, wherein W is oxygen.
- 63. **(Previously presented)** The compound of claim 57, wherein R^{7a} is selected from the group consisting of alkyl, alkenyl, alkynyl, aryl, arylalkyl, and heteroaromatic.
- 64. (Original) The compound of claim 63, wherein R^{7a} is substituted or unsubstituted alkyl.
- 65. (Original) The compound of claim 64, wherein said alkyl is substituted with an aryl group.
- 66. (Original) The compound of claim 63, wherein said substituted or unsubstituted aryl is phenyl.
- 67. (Original) The compound of claim 66, wherein said aryl is substituted with one or more substituents selected from the group consisting of alkyl, alkenyl, alkynyl, aryl, alkoxy, aryloxy,

alkylcarbonyl, arylcarbonyl, alkoxycarbonyl, aryloxycarbonyl, amido, halogen, nitro, azo, alkyl sulfonyl, and arylsulfonyl.

68. (Original) The compound of claim 67, wherein said substituent is alkyl, alkoxy, or nitro.

Claims 69.-81.(Cancelled)

82. (Currently Amended) A pharmaceutical composition comprising a therapeutically effective amount of a substituted tetracycline compound and a pharmaceutically acceptable carrier, wherein said substituted tetracycline is of the formula:

wherein:

X is CHC($R^{13}Y'Y$), $CR^{6'}R^{6}$, S, NR^{6} , or O;

R² is hydrogen, alkyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic or a prodrug moiety;

R⁴ and R⁴ are each hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic or a prodrug moiety;

R², R³, R¹⁰, R¹¹ and R¹² are each hydrogen or a pro-drug moiety;

R⁵ is hydrogen, hydroxyl, or a prodrug moiety;

R⁶, R⁶, and R⁸ are each independently hydrogen, alkyl, alkenyl, alkynyl, aryl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, or halogen;

R⁷ is hydrogen, dialkylamino, or NR^{7c}C(=W')WR^{7a};

R⁸ is hydrogen;

R¹³ is hydrogen, hydroxy, alkyl; alkenyl; alkynyl; alkoxy; alkylthio; alkylsulfinyl; alkylsulfonyl; alkylamino; or an arylalkyl;

Y' and Y are each independently hydrogen; halogen; hydroxyl; cyano, sulfhydryl; amino; alkyl; alkenyl; alkynyl; alkoxy; alkylthio; alkylsulfinyl; alkylsulfonyl; alkylamino; or an arylalkyl;

R⁹ is hydrogen, or NR^{9c}C(=Z')ZR^{9a}; Z is O; Z' is O or S;

R^{9a} is unsubstituted C₃-C₁₀ alkyl, substituted alkyl, substituted or unsubstituted alkenyl, substituted or unsubstituted alkynyl, substituted or unsubstituted alkoxy, substituted or unsubstituted alkylthio, substituted or unsubstituted alkylsulfinyl, substituted or unsubstituted alkylsulfonyl, substituted or unsubstituted arylsulfonyl, substituted or unsubstituted alkoxycarbonyl, substituted or unsubstituted arylcarbonyl, substituted or unsubstituted alkylamino, substituted or unsubstituted arylalkyl, substituted or unsubstituted aryl, substituted or unsubstituted heterocyclic, substituted or unsubstituted heteroaromatic, a steroid, absent, or a prodrug moiety, wherein said substituted alkyl is substituted with halogen, amino, hydroxyl, alkoxy, alkylcarbonyloxy, alkyloxycarbonyl, arylcarbonyloxy, alkoxycarbonylamino, alkoxycarbonyloxy, aryloxycarbonyloxy, carboxylate, alkylcarbonyl, alkylaminoacarbonyl, arylalkyl aminocarbonyl, alkenylaminocarbonyl, alkylcarbonyl, arylcarbonyl, aminoalkyl, arylalkylcarbonyl, alkenylcarbonyl, alkoxycarbonyl, silyl, aminocarbonyl, alkylthiocarbonyl, phosphate, aralkyl, phosphonato, phosphinato, cyano, acylamino, amido, imino, sulfhydryl, alkylthio, sulfate, arylthio, thiocarboxylate, alkylsulfinyl, sulfonato, sulfamoyl, sulfonamido, nitro, cyano, azido, heterocyclyl, alkylaryl, aryl or heteroaryl; further wherein said substituted alkenyl, substituted alkynyl, substituted alkoxy, substituted alkylthio, substituted alkylsulfinyl, substituted alkylsulfonyl, substituted arylsulfonyl, substituted alkoxycarbonyl, substituted arylcarbonyl, substituted alkylamino, substituted arylalkyl, substituted aryl, substituted heterocyclic, or substituted heteroaromatic is substituted with halogen, amino, alkyl, alkenyl, alkynyl, hydroxyl, alkoxy, alkylcarbonyloxy, alkyloxycarbonyl, arylcarbonyloxy, alkoxycarbonylamino, alkoxycarbonyloxy, aryloxycarbonyloxy, carboxylate, alkylcarbonyl, alkylaminoacarbonyl, arylalkyl aminocarbonyl, alkenylaminocarbonyl, alkylcarbonyl, arylcarbonyl, aminoalkyl, arylalkylcarbonyl, alkenylcarbonyl, alkoxycarbonyl, silyl, aminocarbonyl, alkylthiocarbonyl, phosphate, aralkyl, phosphonato, phosphinato, cyano, acylamino, amido, imino, sulfhydryl, alkylthio, sulfate, arylthio, thiocarboxylate, alkylsulfinyl, sulfonato, sulfamoyl, sulfonamido, nitro, cyano, azido, heterocyclyl, alkylaryl, aryl or heteroaryl;

R^{9c} is hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, arylsulfonyl, arylcarbonyl, arylcarbonyl, alkylamino, arylalkyl, aryl, heterocyclic or heteroaromatic;

W is CR^{7d}R^{7e}, NR^{7b} or O;

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W' is O or S; and

R^{7a}, R^{7b}, R^{7c}, R^{7d}, and R^{7e} are each independently hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, arylsulfonyl, alkoxycarbonyl, arylcarbonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic, absent, or a prodrug moiety, and R^{7d} and R^{7e} may be linked to form a ring;

and pharmaceutically acceptable salts thereof, provided that at least one of R9 is not hydrogen when R⁷ is hydrogen or dialkylamino.

83.-102. (Cancelled)

103. (Previously presented) The compound of claim 1, wherein said compound is

and pharmaceutically acceptable salts and prodrugs

thereof.

104. (Previously presented) The compound of claim 1, wherein said compound is

and pharmaceutically acceptable salts and prodrugs

thereof.

105. (Previously presented) The compound of claim 1, wherein said compound is

and pharmaceutically acceptable salts and prodrugs

thereof.

106. (Previously presented) The compound of claim 1, wherein said compound is

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and pharmaceutically acceptable salts and prodrugs

thereof.

(Previously presented) The compound of claim 1, wherein said compound is 107.

and pharmaceutically acceptable salts and prodrugs thereof.

108. (Previously presented) A substituted tetracycline compound, wherein said compound is selected from the group consisting of

and pharmaceutically acceptable salts and prodrugs thereof.

- 109. (Previously presented) The compound of claim 1, wherein said compound is doxycycline 9-carbamic acid 9H-fluoren-9-yl methyl ester and pharmaceutically acceptable salts and prodrugs thereof.
- 110. (Previously presented) The compound of claim 1, wherein said compound is FMOC 9-amino doxycycline and pharmaceutically acceptable salts and prodrugs thereof.

111. **(Previously presented)** The compound of claim 1, wherein said compound is 9-(4'-fluorophenyl) doxycycline carbamate and pharmaceutically acceptable salts and prodrugs thereof.

- 112. **(Previously presented)** The compound of claim 1, wherein said compound is 9-(4'-methoxyphenyl) doxycycline carbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 113. (Currently Amended) The compound of claim 140, wherein said compound is minocycline 9-carbamic acid 9*H*-fluoren-9-yl methyl ester and pharmaceutically acceptable salts and prodrugs thereof.
- 114. (Currently Amended) The compound of claim 140, wherein said compound is FMOC 9-amino minocycline and pharmaceutically acceptable salts and prodrugs thereof.
- 115. (Currently Amended) The compound of claim 140, wherein said compound is 9-(4'-fluorophenyl) minocycline carbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 116. (Currently Amended) The compound of claim 140, wherein said compound is 9-(4'-Methoxyphenyl) minocycline carbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 117. **(Previously presented)** The compound of claim 1, wherein said compound is 9-(2'-bromoethyl) doxycycline carbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 118. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(4'-methylphenyl) sancycline carbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 119. **(Previously presented)** The compound of claim 1, wherein said compound is doxycycline 7-carbamic acid 7*H*-fluoren-7-yl methyl ester and pharmaceutically acceptable salts and prodrugs thereof.
- 120. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(naphthyn-1-yl) doxycycline urea and pharmaceutically acceptable salts and prodrugs thereof.

- 121. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(3-methyl-1-butyl) doxycycline urea and pharmaceutically acceptable salts and prodrugs thereof.
- 122. **(Previously presented)** The compound of claim 1, wherein said compound is 7-phenyl doxycycline urea and pharmaceutically acceptable salts and prodrugs thereof.
- 123. **(Previously presented)** The compound of claim 1, wherein said compound is 7-t-butyl doxycycline urea and pharmaceutically acceptable salts and prodrugs thereof.
- 124. **(Previously presented)** The compound of claim 1, wherein said compound is 7-Fmoc amino doxycycline and pharmaceutically acceptable salts and prodrugs thereof.
- 125. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(4'-chloro-2-trifluoromethylphenyl) doxycycline urea and pharmaceutically acceptable salts and prodrugs thereof.
- 126. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(4'-fluorophenyl) doxycycline carbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 127. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(4'-methoxyphenyl) doxycycline carbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 128. **(Previously presented)** The compound of claim 1, wherein said compound is 7-BOC amino doxycycline and pharmaceutically acceptable salts and prodrugs thereof.
- 129. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(naphthyn-1-yl) doxycycline thiourea 5-propanoic acid ester and pharmaceutically acceptable salts and prodrugs thereof.
- 130. **(Previously presented)** The compound of claim 1, wherein said compound is doxycycline 7-thiocarbamic acid 7*H*-fluoren-7-yl methyl ester and pharmaceutically acceptable salts and prodrugs thereof.

131. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(naphthyn-1-yl) doxycycline thiourea and pharmaceutically acceptable salts and prodrugs thereof.

- 132. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(3-methyl-1-butyl) doxycycline thiourea and pharmaceutically acceptable salts and prodrugs thereof.
- 133. **(Previously presented)** The compound of claim 1, wherein said compound is 7-phenyl amino doxycycline thiourea and pharmaceutically acceptable salts and prodrugs thereof.
- 134. **(Previously presented)** The compound of claim 1, wherein said compound is 7-t-butyl amino doxycycline thiourea and pharmaceutically acceptable salts and prodrugs thereof.
- 135. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(4'-chloro-2'-trifluoromethylphenyl) doxycycline thiourea and pharmaceutically acceptable salts and prodrugs thereof.
- 136. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(4?-fluorophenyl) doxycycline thiocarbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 137. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(4'-methoxyphenyl) doxycycline thiocarbamate and pharmaceutically acceptable salts and prodrugs thereof.
- 138. **(Previously presented)** The compound of claim 1, wherein said compound is 7-(naphthyn-1-yl) doxycycline urea 5-propanoic acid ester and pharmaceutically acceptable salts and prodrugs thereof.
- 139. **(Previously presented)** A substituted tetracycline compound, wherein said compound is selected from the group consisting of:

9-neopentyl minocycline carbamate;

9-BOC amino doxycycline;

9-(n-pentyl) minocycline carbamate;

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9-BOC amino minocycline carbamate;

9-(n-pentyl) minocycline carbamate;

9-prop-2'-enyl minocycline carbamate;

9-ethyl minocycline carbamate;

9-n-butyl minocycline carbamate

9-n-but-3-enyl minocycline carbamate; and

9-i-butyl minocycline carbamate; and pharmaceutically acceptable salts and prodrugs thereof.

140. (New) A substituted tetracycline compound, wherein said compound is of the formula:

wherein:

X is CHC(R¹³Y'Y), CR⁶'R⁶, S, NR⁶, or O;

R² is hydrogen, alkyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic or a prodrug moiety;

R⁴ and R⁴ are each hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic or a prodrug moiety;

R²', R³, R¹⁰, R¹¹ and R¹² are each hydrogen or a pro-drug moiety;

R⁵ is hydrogen, hydroxyl, or a prodrug moiety;

R⁶, R⁶, and R⁸ are each independently hydrogen, alkyl, alkenyl, alkynyl, aryl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, or halogen;

R⁷ dialkylamino;

R⁸ is hydrogen;

R¹³ is hydrogen, hydroxy, alkyl; alkenyl; alkynyl; alkoxy; alkylthio; alkylsulfinyl; alkylsulfonyl; alkylamino; or an arylalkyl;

Y' and Y are each independently hydrogen; halogen; hydroxyl; cyano, sulfhydryl; amino; alkyl; alkenyl; alkoxy; alkylthio; alkylsulfinyl; alkylsulfonyl; alkylamino; or an arylalkyl;

$$R^9$$
 is $NR^{9c}C(=Z')ZR^{9a}$;

Z is O;

Z' is O or S;

R^{9a} is unsubstituted or substituted alkyl, substituted or unsubstituted alkenyl, substituted or unsubstituted alkynyl, substituted or unsubstituted alkoxy, substituted or unsubstituted alkylthio, substituted or unsubstituted alkylsulfinyl, substituted or unsubstituted alkylsulfonyl, substituted or unsubstituted arylsulfonyl, substituted or unsubstituted alkoxycarbonyl, substituted or unsubstituted arylcarbonyl, substituted or unsubstituted alkylamino, substituted or unsubstituted arylalkyl, substituted or unsubstituted aryl, substituted or unsubstituted heterocyclic, substituted or unsubstituted heteroaromatic, absent, or a prodrug moiety, wherein said substituted alkyl is substituted with halogen, hydroxyl, alkoxy, alkylcarbonyloxy, alkyloxycarbonyl, arylcarbonyloxy, alkoxycarbonylamino, alkoxycarbonyloxy, aryloxycarbonyloxy, carboxylate, alkylcarbonyl, alkylaminoacarbonyl, arylalkyl aminocarbonyl, alkenylaminocarbonyl, alkylcarbonyl, arylcarbonyl, aminoalkyl, arylalkylcarbonyl, alkenylcarbonyl, alkoxycarbonyl, silyl, aminocarbonyl, alkylthiocarbonyl, phosphate, aralkyl, phosphonato, phosphinato, cyano, acylamino, amido, imino, sulfhydryl, alkylthio, sulfate, arylthio, thiocarboxylate, alkylsulfinyl, sulfonato, sulfamoyl, sulfonamido, nitro, cyano, azido, heterocyclyl, alkylaryl, aryl or heteroaryl; further wherein said substituted alkenyl, substituted alkynyl, substituted alkoxy, substituted alkylthio, substituted alkylsulfinyl, substituted alkylsulfonyl, substituted arylsulfonyl, substituted alkoxycarbonyl, substituted arylcarbonyl, substituted alkylamino, substituted arylalkyl, substituted aryl, substituted heterocyclic, or substituted heteroaromatic is substituted with halogen, amino, alkyl, alkenyl, alkynyl, hydroxyl, alkoxy, alkylcarbonyloxy, alkyloxycarbonyl, arylcarbonyloxy, alkoxycarbonylamino, alkoxycarbonyloxy, aryloxycarbonyloxy, carboxylate, alkylcarbonyl, alkylaminoacarbonyl, arylalkyl aminocarbonyl, alkenylaminocarbonyl, alkylcarbonyl, arylcarbonyl, aminoalkyl, arylalkylcarbonyl, alkenylcarbonyl, alkoxycarbonyl, silyl, aminocarbonyl, alkylthiocarbonyl, phosphate, aralkyl, phosphonato, phosphinato, cyano, acylamino, amido, imino, sulfhydryl, alkylthio, sulfate, arylthio, thiocarboxylate, alkylsulfinyl, sulfonato, sulfamoyl, sulfonamido, nitro, cyano, azido, heterocyclyl, alkylaryl, aryl or heteroaryl; and

R^{9c} is hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, arylsulfonyl, alkoxycarbonyl, arylcarbonyl, alkylamino, arylalkyl, aryl, heterocyclic or heteroaromatic; and pharmaceutically acceptable salts thereof.

141. (New) The compound of claim 140, wherein R², R², R³, R⁸, R¹⁰, R¹¹, and R¹² are each hydrogen.

142. (New) The compound of claim 140, wherein R⁴ and R⁴ are each methyl

- 143. (New) The compound of claim 140, wherein R⁵ is hydrogen.
- 144. (New) The compound of claim 140, wherein R⁹ is NR^{9c}C(=Z')ZR^{9a}.
- 145. (New) The compound of claim 144, wherein R^{9c} is hydrogen.
- 146. (New) The compound of claim 144, wherein Z' is oxygen.
- 147. (New) The compound of claim 144, wherein Z' is sulfur.
- 148. (New) The compound of claim 144, wherein Z is oxygen.
- 149. (New) The compound of claim 144, wherein R^{9a} is substituted alkyl, alkynyl, aryl, arylalkyl, or heteroaromatic.
- 150. **(New)** The compound of claim 149, wherein said substituted alkyl is substituted with one or more substituents selected from the group consisting of alkoxycarbonyl, arylcarbonyl, halogen, hydroxy, alkylamino, alkoxy, or aryl.
- 151. (New) The compound of claim 149, wherein R^{9a} is substituted or unsubstituted aryl.
- 152. (New) The compound of claim 151, wherein said substituted or unsubstituted aryl is phenyl.
- 153. (New) The compound of claim 151, wherein said aryl is substituted with one or more substituents selected from the group consisting of alkyl, alkenyl, alkynyl, aryl, alkoxy, aryloxy, alkylcarbonyl, arylcarbonyl, alkoxycarbonyl, aryloxycarbonyl, amido, halogen, nitro, azo, alkyl sulfonyl, and arylsulfonyl.

154. (New) A pharmaceutical composition comprising a therapeutically effective amount of a substituted tetracycline compound and a pharmaceutically acceptable carrier, wherein said substituted tetracycline is of the formula:

$$R^{0}$$

$$R^{0$$

(I)

wherein:

X is CHC(R¹³Y'Y), CR⁶'R⁶, S, NR⁶, or O;

R² is hydrogen, alkyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic or a prodrug moiety;

R⁴ and R⁴ are each hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, aryl, heterocyclic, heteroaromatic or a prodrug moiety;

R², R³, R¹⁰, R¹¹ and R¹² are each hydrogen or a pro-drug moiety;

R⁵ is hydrogen, hydroxyl, or a prodrug moiety;

R⁶, R⁶, and R⁸ are each independently hydrogen, alkyl, alkenyl, alkynyl, aryl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, alkylamino, arylalkyl, or halogen;

R⁷ dialkylamino;

R⁸ is hydrogen;

R¹³ is hydrogen, hydroxy, alkyl; alkenyl; alkynyl; alkoxy; alkylthio; alkylsulfinyl; alkylsulfonyl; alkylamino; or an arylalkyl;

Y' and Y are each independently hydrogen; halogen; hydroxyl; cyano, sulfhydryl; amino; alkyl; alkenyl; alkoxy; alkylthio; alkylsulfinyl; alkylsulfonyl; alkylamino; or an arylalkyl;

 R^9 is $NR^{9c}C(=Z')ZR^{9a}$;

Z is O;

Z' is O or S;

R^{9a} is unsubstituted or substituted alkyl, substituted or unsubstituted alkenyl, substituted or unsubstituted or unsubstituted alkynyl, substituted or unsubstituted alkylsulfinyl, substituted or unsubstituted alkylsulfinyl, substituted or unsubstituted alkylsulfonyl, substituted or unsubstituted arylsulfonyl, substituted or unsubstituted or u

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unsubstituted arylcarbonyl, substituted or unsubstituted alkylamino, substituted or unsubstituted arylalkyl, substituted or unsubstituted aryl, substituted or unsubstituted heterocyclic, substituted or unsubstituted heteroaromatic, absent, or a prodrug moiety, wherein said substituted alkyl is substituted with halogen, hydroxyl, alkoxy, alkylcarbonyloxy, alkyloxycarbonyl, arylcarbonyloxy, alkoxycarbonylamino, alkoxycarbonyloxy, aryloxycarbonyloxy, carboxylate, alkylcarbonyl, alkylaminoacarbonyl, arylalkyl aminocarbonyl, alkenylaminocarbonyl, alkylcarbonyl, arylcarbonyl, aminoalkyl, arylalkylcarbonyl, alkenylcarbonyl, alkoxycarbonyl, silyl, aminocarbonyl, alkylthiocarbonyl, phosphate, aralkyl, phosphonato, phosphinato, cyano, acylamino, amido, imino, sulfhydryl, alkylthio, sulfate, arylthio, thiocarboxylate, alkylsulfinyl, sulfonato, sulfamoyl, sulfonamido, nitro, cyano, azido, heterocyclyl, alkylaryl, aryl or heteroaryl; further wherein said substituted alkenyl, substituted alkynyl, substituted alkoxy, substituted alkylthio, substituted alkylsulfinyl, substituted alkylsulfonyl, substituted arylsulfonyl, substituted alkoxycarbonyl, substituted arylcarbonyl, substituted alkylamino, substituted arylalkyl, substituted aryl, substituted heterocyclic, or substituted heteroaromatic is substituted with halogen, amino, alkyl, alkenyl, alkynyl, hydroxyl, alkoxy, alkylcarbonyloxy, alkyloxycarbonyl, arylcarbonyloxy, alkoxycarbonylamino, alkoxycarbonyloxy, aryloxycarbonyloxy, carboxylate, alkylcarbonyl, alkylaminoacarbonyl, arylalkyl aminocarbonyl, alkenylaminocarbonyl, alkylcarbonyl, arylcarbonyl, aminoalkyl, arylalkylcarbonyl, alkenylcarbonyl, alkoxycarbonyl, silyl, aminocarbonyl, alkylthiocarbonyl, phosphate, aralkyl, phosphonato, phosphinato, cyano, acylamino, amido, imino, sulfhydryl, alkylthio, sulfate, arylthio, thiocarboxylate, alkylsulfinyl, sulfonato, sulfamoyl, sulfonamido, nitro, cyano, azido, heterocyclyl, alkylaryl, aryl or heteroaryl; and

R^{9c} is hydrogen, alkyl, alkenyl, alkynyl, alkoxy, alkylthio, alkylsulfinyl, alkylsulfonyl, arylsulfonyl, alkoxycarbonyl, arylcarbonyl, alkylamino, arylalkyl, aryl, heterocyclic or heteroaromatic; and pharmaceutically acceptable salts thereof.